

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

TERMINAL (ENTER 1, 2, 3, OR ?):2

\* \* \* \* \* Welcome to STN International \* \* \* \* \*

NEWS 1 Web Page URLs for STN Seminar Schedule - N. America  
NEWS 2 "Ask CAS" for self-help around the clock  
NEWS 3 FEB 28 PATDPAFULL - New display fields provide for legal status  
data from INPADOC  
NEWS 4 FEB 28 BABS - Current-awareness alerts (SDIs) available  
NEWS 5 MAR 02 GBFULL: New full-text patent database on STN  
NEWS 6 MAR 03 REGISTRY/ZREGISTRY - Sequence annotations enhanced  
NEWS 7 MAR 03 MEDLINE file segment of TOXCENTER reloaded  
NEWS 8 MAR 22 KOREAPAT now updated monthly; patent information enhanced  
NEWS 9 MAR 22 Original IDE display format returns to REGISTRY/ZREGISTRY  
NEWS 10 MAR 22 PATDPASPC - New patent database available  
NEWS 11 MAR 22 REGISTRY/ZREGISTRY enhanced with experimental property tags  
NEWS 12 APR 04 EPFULL enhanced with additional patent information and new  
fields  
NEWS 13 APR 04 EMBASE - Database reloaded and enhanced  
NEWS 14 APR 18 New CAS Information Use Policies available online  
NEWS 15 APR 25 Patent searching, including current-awareness alerts (SDIs),  
based on application date in CA/Caplus and USPATFULL/USPAT2  
may be affected by a change in filing date for U.S.  
applications.  
NEWS 16 APR 28 Improved searching of U.S. Patent Classifications for  
U.S. patent records in CA/Caplus  
NEWS 17 MAY 23 GBFULL enhanced with patent drawing images  
NEWS 18 MAY 23 REGISTRY has been enhanced with source information from  
CHEMCATS  
NEWS 19 JUN 06 The Analysis Edition of STN Express with Discover!  
(Version 8.0 for Windows) now available  
NEWS 20 JUN 13 RUSSIAPAT: New full-text patent database on STN  
NEWS 21 JUN 13 FRFULL enhanced with patent drawing images  
NEWS 22 JUN 27 MARPAT displays enhanced with expanded G-group definitions  
and text labels  
NEWS 23 JUL 01 MEDICONF removed from STN  
NEWS 24 JUL 07 STN Patent Forums to be held in July 2005  
NEWS 25 JUL 13 SCISEARCH reloaded

NEWS EXPRESS JUNE 13 CURRENT WINDOWS VERSION IS V8.0, CURRENT  
MACINTOSH VERSION IS V6.0c(ENG) AND V6.0Jc(JP),  
AND CURRENT DISCOVER FILE IS DATED 13 JUNE 2005

NEWS HOURS STN Operating Hours Plus Help Desk Availability  
NEWS INTER General Internet Information  
NEWS LOGIN Welcome Banner and News Items  
NEWS PHONE Direct Dial and Telecommunication Network Access to STN  
NEWS WWW CAS World Wide Web Site (general information)

Enter NEWS followed by the item number or name to see news on that  
specific topic.

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agreement. Please note that this agreement limits use to scientific research. Use for software development or design or implementation of commercial gateways or other similar uses is prohibited and may result in loss of user privileges and other penalties.

\* \* \* \* \* STN Columbus \* \* \* \* \*

FILE 'HOME' ENTERED AT 06:25:36 ON 19 JUL 2005

=> file reg

COST IN U.S. DOLLARS

SINCE FILE

ENTRY

TOTAL

SESSION

FULL ESTIMATED COST

0.84

0.84

FILE 'REGISTRY' ENTERED AT 06:27:51 ON 19 JUL 2005

USE IS SUBJECT TO THE TERMS OF YOUR STN CUSTOMER AGREEMENT.

PLEASE SEE "HELP USAGETERMS" FOR DETAILS.

COPYRIGHT (C) 2005 American Chemical Society (ACS)

Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

DICTIONARY FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

New CAS Information Use Policies, enter HELP USAGETERMS for details.

TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

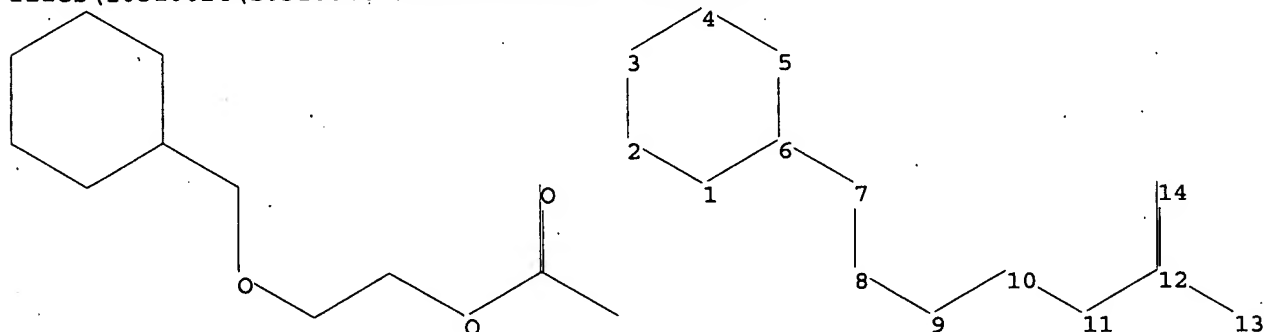
\*\*\*\*\*  
\*  
\* The CA roles and document type information have been removed from \*  
\* the IDE default display format and the ED field has been added, \*  
\* effective March 20, 2005. A new display format, IDERL, is now \*  
\* available and contains the CA role and document type information. \*  
\*  
\*\*\*\*\*

Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

=>

Uploading C:\Documents and Settings\PZucker\My Documents\Examination Auxillary files\10510024\10510024 unsubstituted core.str

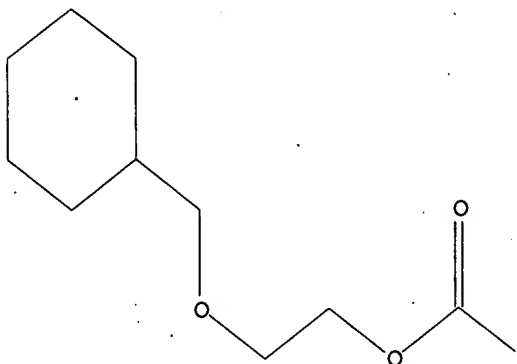


chain nodes :  
 7 8 9 10 11 12 13 14  
 ring nodes :  
 1 2 3 4 5 6  
 chain bonds :  
 6-7 7-8 8-9 9-10 10-11 11-12 12-13 12-14  
 ring bonds :  
 1-2 1-6 2-3 3-4 4-5 5-6  
 exact/norm bonds :  
 1-2 1-6 2-3 3-4 4-5 5-6 7-8 8-9 10-11 11-12 12-14  
 exact bonds :  
 6-7 9-10 12-13

Match level :  
 1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS  
 11:CLASS 12:CLASS 13:CLASS 14:CLASS

L1 STRUCTURE UPLOADED

=> d l1  
 L1 HAS NO ANSWERS  
 L1 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l1 sss sam  
 SAMPLE SEARCH INITIATED 06:28:12 FILE 'REGISTRY'  
 SAMPLE SCREEN SEARCH COMPLETED - 14803 TO ITERATE

13.5% PROCESSED 2000 ITERATIONS  
 INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)  
 SEARCH TIME: 00.00.01

50 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
 BATCH \*\*COMPLETE\*\*  
 PROJECTED ITERATIONS: 288773 TO 303347  
 PROJECTED ANSWERS: 8443 TO 11095

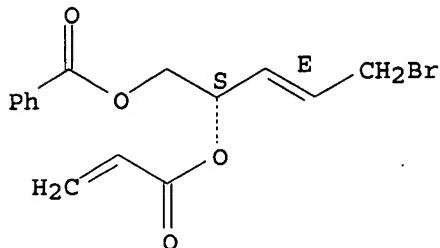
L2 50 SEA SSS SAM L1

=> d scan

L2 50 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 2-Propenoic acid, (1S,2E)-1-[(benzyloxy)methyl]-4-bromo-2-butenyl ester

(9CI)  
MF C15 H15 Br O4

Absolute stereochemistry. Rotation (+).  
Double bond geometry as shown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> d cost

COST IN U.S. DOLLARS

CONNECT CHARGES

NETWORK CHARGES

FULL ESTIMATED COST

SINCE FILE

ENTRY

0.74

0.12

0.86

TOTAL

SESSION

1.34

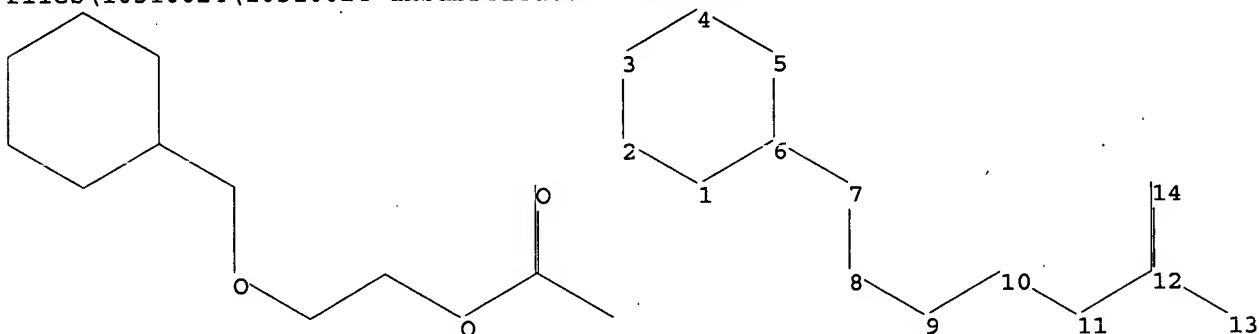
0.36

1.70

IN FILE 'REGISTRY' AT 06:28:49 ON 19 JUL 2005

=>

Uploading C:\Documents and Settings\PZucker\My Documents\Examination Auxillary  
files\10510024\10510024 unsubstituted core2.str



chain nodes :

7 8 9 10 11 12 13 14

ring nodes :

1 2 3 4 5 6

chain bonds :

6-7 7-8 8-9 9-10 10-11 11-12 12-13 12-14

ring bonds :

1-2 1-6 2-3 3-4 4-5 5-6

exact/norm bonds :

1-2 1-6 2-3 3-4 4-5 5-6 7-8 8-9 10-11 11-12 12-14

exact bonds :

6-7 9-10 12-13

Hydrogen count :

1:>= minimum 2 2:>= minimum 2 3:>= minimum 2 5:>= minimum 2 6:>= minimum 1

Match level :

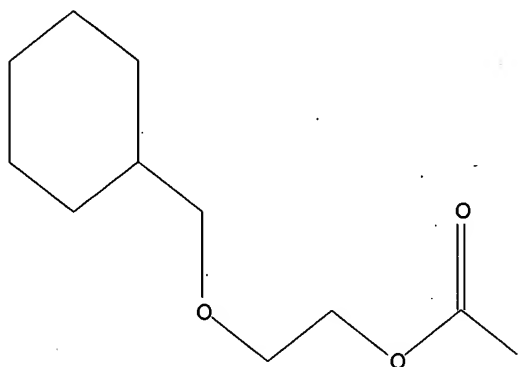
1:Atom 2:Atom 3:Atom 4:Atom 5:Atom 6:Atom 7:CLASS 8:CLASS 9:CLASS 10:CLASS  
11:CLASS 12:CLASS 13:CLASS 14:CLASS

L3 STRUCTURE UPLOADED

=> d l3

L3 HAS NO ANSWERS

L3 STR



Structure attributes must be viewed using STN Express query preparation.

=> search l3 sss sam

SAMPLE SEARCH INITIATED 06:30:36 FILE 'REGISTRY'

SAMPLE SCREEN SEARCH COMPLETED - 14803 TO ITERATE

13.5% PROCESSED 2000 ITERATIONS  
INCOMPLETE SEARCH (SYSTEM LIMIT EXCEEDED)  
SEARCH TIME: 00.00.01

0 ANSWERS

FULL FILE PROJECTIONS: ONLINE \*\*COMPLETE\*\*  
BATCH \*\*COMPLETE\*\*

PROJECTED ITERATIONS: 288773 TO 303347  
PROJECTED ANSWERS: 0 TO 0

L4 0 SEA SSS SAM L3

=> search.l3 sss full

FULL SEARCH INITIATED 06:31:36 FILE 'REGISTRY'

FULL SCREEN SEARCH COMPLETED - 296742 TO ITERATE

100.0% PROCESSED 296742 ITERATIONS  
SEARCH TIME: 00.00.02

74 ANSWERS

L5 74 SEA SSS FUL L3

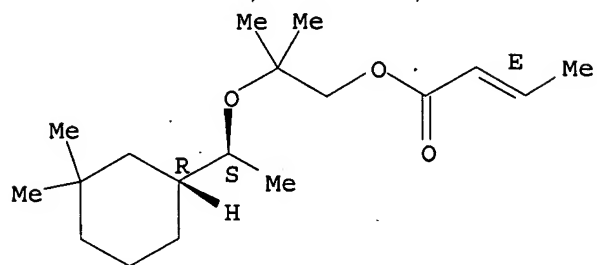
=> d scan

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN

IN 2-Butenoic acid, 2-[(1S)-1-[(1R)-3,3-dimethylcyclohexyl]ethoxy]-2-methylpropyl ester, (2E)- (9CI)

MF C18 H32 O3

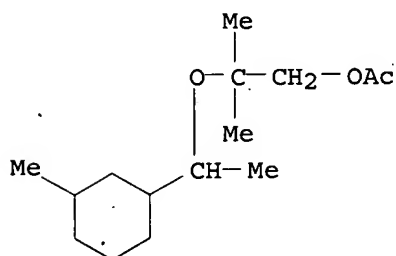
Absolute stereochemistry.  
Double bond geometry as shown.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

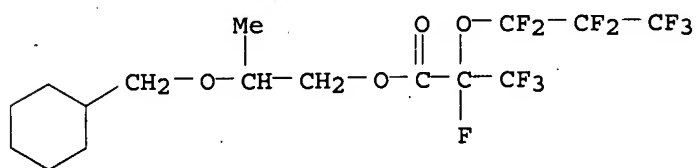
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN 1-Propanol, 2-methyl-2-[1-(3-methylcyclohexyl)ethoxy]-, acetate (9CI)  
MF C15 H28 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

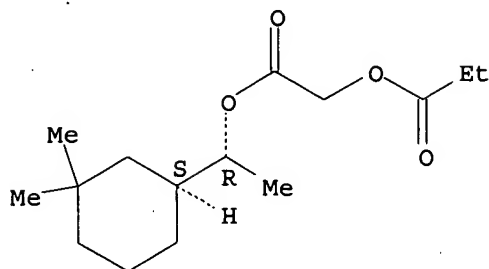
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN Propanoic acid, 2,3,3,3-tetrafluoro-2-(heptafluoropropoxy)-,  
2-(cyclohexylmethoxy)propyl ester (9CI)  
MF C16 H19 F11 O4



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN Acetic acid, (1-oxopropoxy)-, (1R)-1-[(1S)-3,3-dimethylcyclohexyl]ethyl  
ester (9CI)  
MF C15 H26 O4

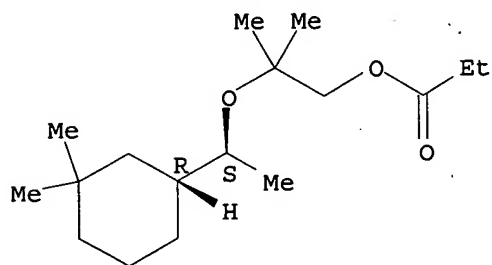
Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 1-Propanol, 2-[1-(3,3-dimethylcyclohexyl)ethoxy]-2-methyl-, propanoate,  
 (R\*,S\*)-(9CI)  
 MF C17 H32 O3

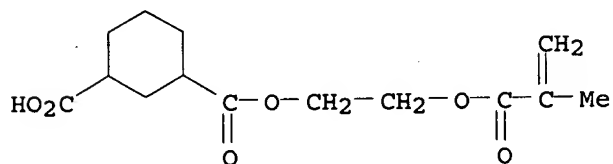
Relative stereochemistry.



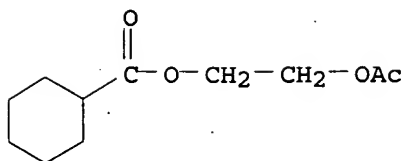
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 1,3-Cyclohexanedicarboxylic acid, mono[2-[(2-methyl-1-oxo-2-propenyl)oxy]ethyl] ester, homopolymer (9CI)  
 MF (C14 H20 O6)x  
 CI PMS

CM 1

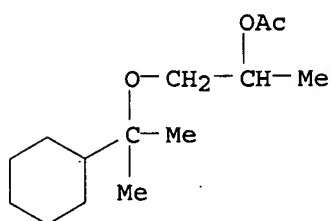


L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Cyclohexanecarboxylic acid, 2-hydroxyethyl ester, acetate (8CI)  
 MF C11 H18 O4



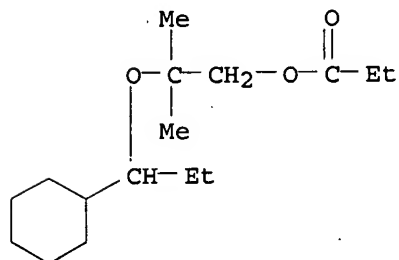
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 2-Propanol, 1-(1-cyclohexyl-1-methylethoxy)-, acetate (9CI)  
 MF C14 H26 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

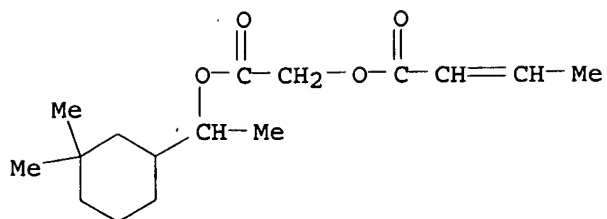
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN.  
 IN 1-Propanol, 2-(1-cyclohexylpropoxy)-2-methyl-, propanoate (9CI)  
 MF C16 H30 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

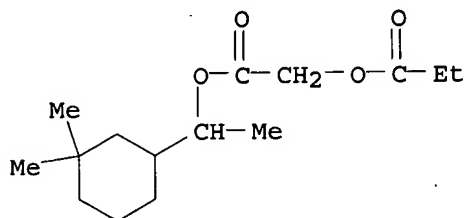
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 2-Butenoic acid, 2-[1-(3,3-dimethylcyclohexyl)ethoxy]-2-oxoethyl ester  
 (9CI)  
 MF C16 H26 O4





\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

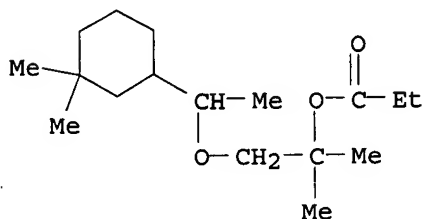
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Acetic acid, (1-oxopropoxy)-, 1-(3,3-dimethylcyclohexyl)ethyl ester (9CI)  
 MF C15 H26 O4



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

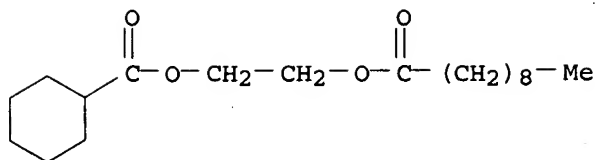
HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 2-Propanol, 1-[1-(3,3-dimethylcyclohexyl)ethoxy]-2-methyl-, propanoate (9CI)  
 MF C17 H32 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

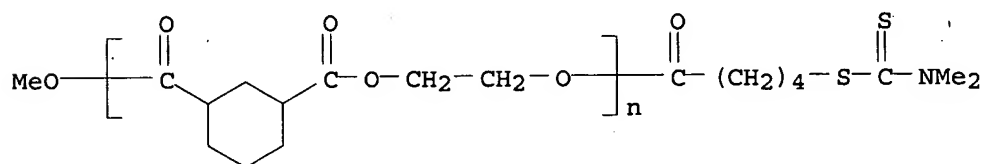
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Decanoic acid, 2-hydroxyethyl ester, cyclohexanecarboxylate (8CI)  
 MF C19 H34 O4



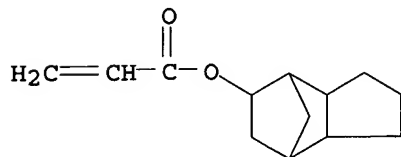
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 2-Propenoic acid, 2-hydroxyethyl ester, polymer with  $\alpha$ -[5-  
 [[(dimethylamino)thioxomethyl]thio]-1-oxopentyl]- $\omega$ -methoxypoly(oxy-  
 1,2-ethanediylloxycarbonyl-1,3-cyclohexanediylcarbonyl) and  
 octahydro-4,7-methano-1H-inden-5-yl 2-propenoate, diblock (9CI)  
 MF (C13 H18 O2 . (C10 H14 O4)n C9 H17 N O2 S2 . C5 H8 O3)x  
 CI PMS

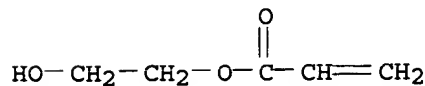
CM 1



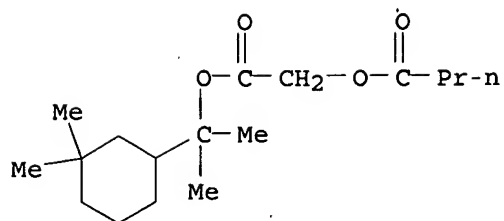
CM 2



CM 3

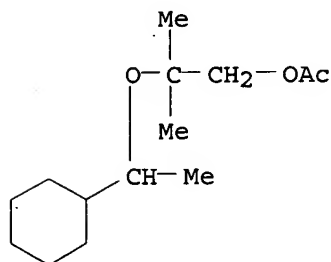


L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Butanoic acid, 2-[1-(3,3-dimethylcyclohexyl)-1-methylethoxy]-2-oxoethyl  
 ester (9CI)  
 MF C17 H30 O4



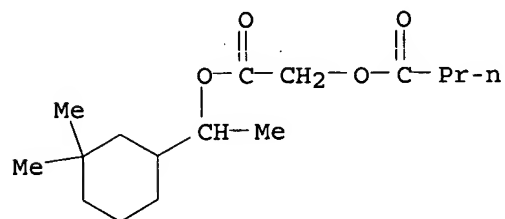
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 1-Propanol, 2-(1-cyclohexylethoxy)-2-methyl-, acetate (9CI)  
 MF C14 H26 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

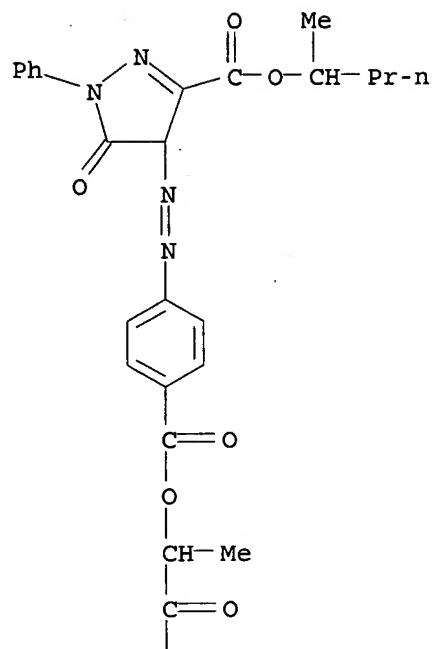
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Butanoic acid, 2-[1-(3,3-dimethylcyclohexyl)ethoxy]-2-oxoethyl ester (9CI)  
 MF C16 H28 O4



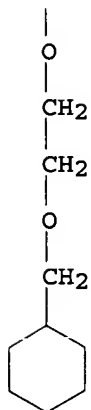
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 1H-Pyrazole-3-carboxylic acid, 4-[[4-[[2-[2-(cyclohexylmethoxy)ethoxy]-1-methyl-2-oxoethoxy]carbonyl]phenyl]azo]-4,5-dihydro-5-oxo-1-phenyl-, 1-methylbutyl ester (9CI)  
 MF C34 H42 N4 O8

PAGE 1-A

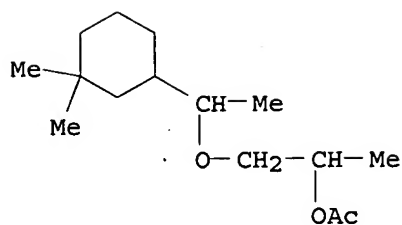


PAGE 2-A



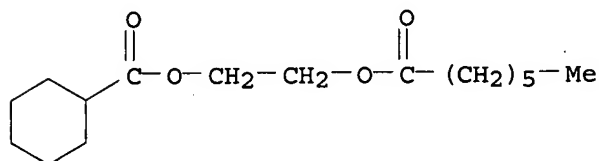
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN 2-Propanol, 1-[1-(3,3-dimethylcyclohexyl)ethoxy]-, acetate (9CI)  
MF C15 H28 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

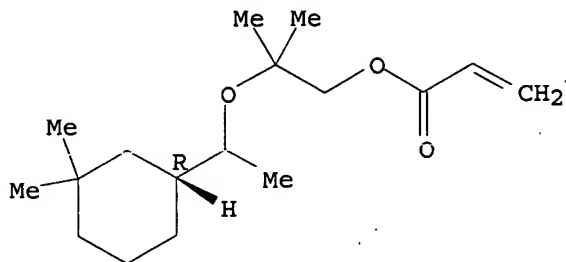
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Heptanoic acid, 2-hydroxyethyl ester cyclohexanecarboxylate (8CI)  
 MF C16 H28 O4



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN 2-Propenoic acid, 2-[1-[(1R)-3,3-dimethylcyclohexyl]ethoxy]-2-methylpropyl  
 ester (9CI)  
 MF C17 H30 O3

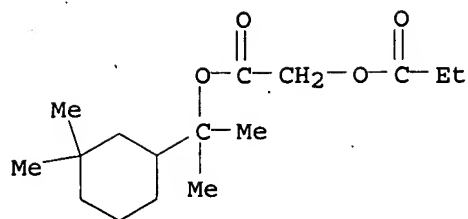
Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):10

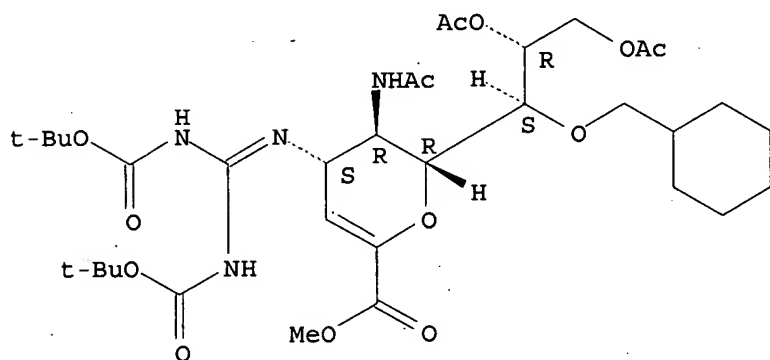
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Acetic acid, (1-oxopropoxy)-, 1-(3,3-dimethylcyclohexyl)-1-methylethyl  
 ester (9CI)  
 MF C16 H28 O4



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN D-glycero-D-galacto-Non-2-enonic acid, 5-(acetylamino)-2,6-anhydro-4-  
 [[bis[[[(1,1-dimethylethoxy) carbonyl] amino] methylene] amino]-7-O-  
 (cyclohexylmethyl)-3,4,5-trideoxy-, methyl ester, 8,9-diacetate (9CI)  
 MF C34 H54 N4 O13

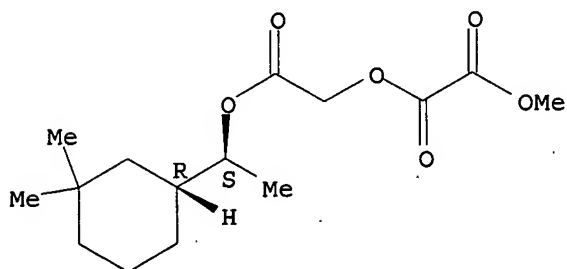
Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
 IN Ethanedioic acid, 2-[(1S)-1-[(1R)-3,3-dimethylcyclohexyl]ethoxy]-2-  
 oxoethyl methyl ester (9CI)  
 MF C15 H24 O6

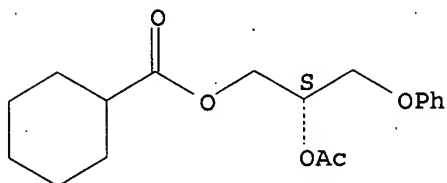
Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

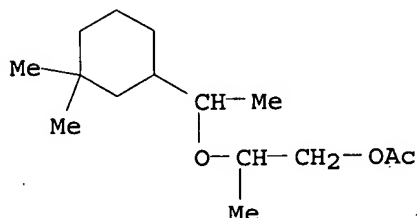
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN Cyclohexanecarboxylic acid, 2-(acetyloxy)-3-phenoxypropyl ester, (S)-  
(9CI)  
MF C18 H24 O5

Absolute stereochemistry.



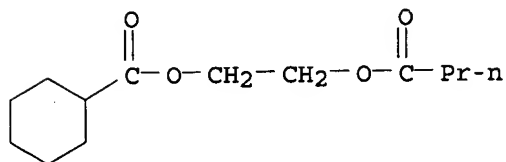
\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN 1-Propanol, 2-[1-(3,3-dimethylcyclohexyl)ethoxy]-, acetate (9CI)  
MF C15 H28 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN Cyclohexanecarboxylic acid, 2-hydroxyethyl ester butyrate (8CI)  
MF C13 H22 O4

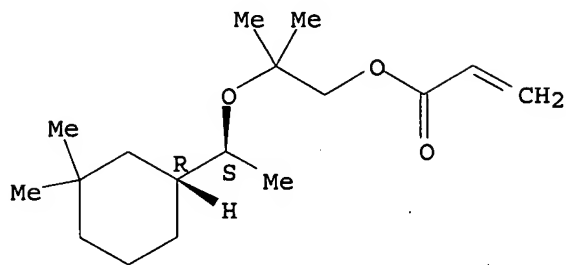


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN 2-Propenoic acid, 2-[(1S)-1-[(1R)-3,3-dimethylcyclohexyl]ethoxy]-2-methylpropyl ester (9CI)

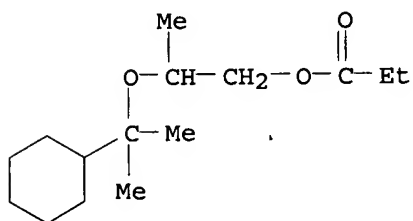
MF C17 H30 O3

Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

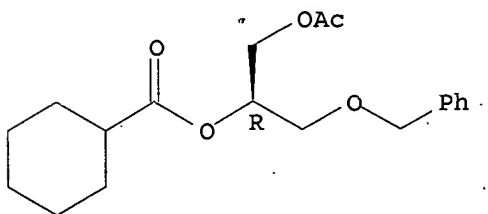
L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN 1-Propanol, 2-(1-cyclohexyl-1-methylethoxy)-, propanoate (9CI)  
MF C15 H28 O3



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN Cyclohexanecarboxylic acid, (1R)-1-[(acetyloxy)methyl]-2-(phenylmethoxy)ethyl ester (9CI)  
MF C19 H26 O5

Absolute stereochemistry. Rotation (-).

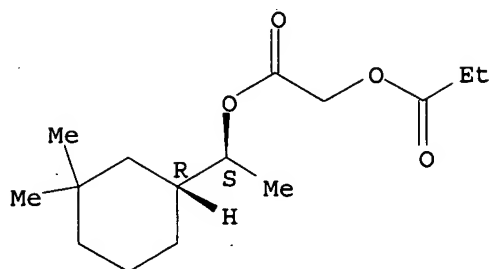


\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

L5 74 ANSWERS REGISTRY COPYRIGHT 2005 ACS on STN  
IN Acetic acid, (1-oxopropoxy)-, (1S)-1-[(1R)-3,3-dimethylcyclohexyl]ethyl ester (9CI)  
MF C15 H26 O4



Absolute stereochemistry.



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

HOW MANY MORE ANSWERS DO YOU WISH TO SCAN? (1):0

=> save temp l5 rawfinds/a  
ANSWER SET L5 HAS BEEN SAVED AS 'RAWFINDS/A'

=> file caplus'  
'CAPLUS'' IS NOT A VALID FILE NAME  
SESSION CONTINUES IN FILE 'REGISTRY'  
Enter "HELP FILE NAMES" at an arrow prompt (=>) for a list of files  
that are available. If you have requested multiple files, you can  
specify a corrected file name or you can enter "IGNORE" to continue  
accessing the remaining file names entered.

=> file caplus		
COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	165.20	166.04

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FILE COVERS 1907 - 19 Jul 2005 VOL 143 ISS 4  
FILE LAST UPDATED: 18 Jul 2005 (20050718/ED)

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substance identification.

=> l5  
L6 21 L5  
=> save temp odorrefs/a

ENTER L#, L# RANGE, ALL, OR (END):16  
ANSWER SET L6 HAS BEEN SAVED AS 'ODORREFS/A'

=> odor

64767 ODOR  
11508 ODORS  
L7 69532 ODOR  
(ODOR OR ODORS)

=> 16 and 17

L8 7 L6 AND L7

=> d 18 1-7 ti

L8 ANSWER 1 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Preparation of alicyclic ester with musk fragrance for perfumes or body care products

L8 ANSWER 2 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Odorous compositions containing (1S,1'R)-2-[1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy]-2-methyl-1-propanol alkenoate esters

L8 ANSWER 3 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Methods for the production of novel alicyclic esters having a musky smell

L8 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Preparation of cycloalkanecarboxylic acid derivatives for use as fragrances with musk characteristics

L8 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI A method for the calculation of odor character from molecular structure

L8 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Preparation of esters with musky odor and their use in perfumery

L8 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Synthesis of ethylene glycol mixed esters of alkanolic and cyclohexanoic acids

=> d 18 1-7 ti fbib abs

L8 ANSWER 1 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Preparation of alicyclic ester with musk fragrance for perfumes or body care products  
AN 2005:120865 CAPLUS  
DN 142:198227  
TI Preparation of alicyclic ester with musk fragrance for perfumes or body care products  
IN Eh, Marcus  
PA Symrise GmbH & Co. KG, Germany  
SO PCT Int. Appl., 42 pp.  
CODEN: PIXXD2  
DT Patent  
LA German  
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2005012222	A1	20050210	WO 2004-EP51361	20040705
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TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW  
 RW: BW, GH, GM, KE, LS, MW, MZ, NA, SD, SL, SZ, TZ, UG, ZM, ZW, AM,  
 AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK,  
 EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PL, PT, RO, SE,  
 SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE,  
 SN, TD, TG

DE 2003-10335053

A 20030731

DE 10335053

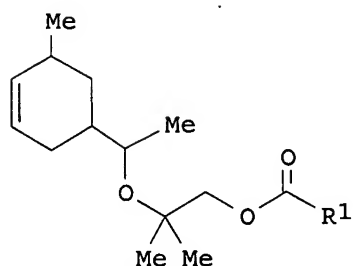
A1 20050224

DE 2003-10335053

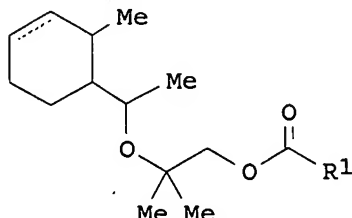
20030731

OS CASREACT 142:198227; MARPAT 142:198227

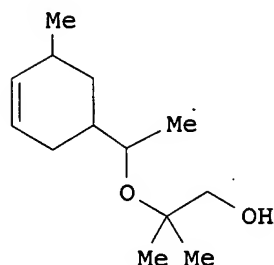
GI



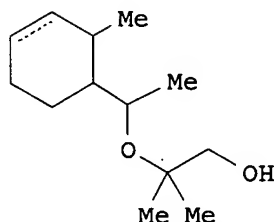
I



II



III



IV

AB A compds. I [R1 = (un)substituted, (un)branched C1-5-alkyl, (un)branched C2-5-alkylene, C3-5-cycloalkyl] and II [dotted line may be either a single or a double bond] are disclosed. The procedure for the preparation of I comprises the esterification with alcs. III and IV. Thus, propionic acid 2-methyl-2-[1-(2-methylcyclohexyl)ethoxy]propyl ester [II; R1 = Et, dashed line = single bond] was prepared from MeCOC6H4Me-2 via hydrogenation with Ru/C, reaction with isobutylene oxidation in the presence of BF3·OEt2, and acylation with (EtCO)2O in the presence of Et3N and catalytic DMAP. A perfume formulation containing I (R1 = Et, dashed line = single bond) was prepared

RE.CNT 2 THERE ARE 2 CITED REFERENCES AVAILABLE FOR THIS RECORD  
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 2 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN

TI Odorous compositions containing (1S,1'R)-2-[1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy]-2-methyl-1-propanol alkenoate esters

AN 2004:780341 CAPLUS

DN 141:301056

TI Odorous compositions containing (1S,1'R)-2-[1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy]-2-methyl-1-propanol alkenoate esters

IN Williams, Alvin Scott

PA Switz.

SO U.S. Pat. Appl. Publ., 6 pp.

CODEN: USXXCO

DT Patent  
LA English  
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	US 2004186043	A1	20040923	US 2004-792375	20040302
				WO 2003-IB1079	A 20030319
	EP 1459735	A1	20040922	EP 2004-100807	20040301
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, PL, SK			WO 2003-IB1079	A 20030319
	JP 2004285357	A2	20041014	JP 2004-81377	20040319
				WO 2003-IB1079	A 20030319

OS MARPAT 141:301056

AB The present invention relates to the field of perfumery, and more precisely to a compound such as (1S,1'R)-2-[1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy]-2-methylpropyl 2-propenoate (I). The present invention concerns also the use of the compound in the perfumery industry as well as the compns. or articles associated with the compound. Thus, to a suspension of NaH (0.368 mol) in 1000 mL of dry THF at 0° were added dropwise 70.0 g (1S,1'R)-2-[1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy]-2-methyl-1-propanol followed by the addition of acryloyl chloride. A 96% pure 2-[1-(3',3'-dimethyl-1'-cyclohexyl)ethoxy]-2-methylpropyl 2-propenoate in the form of a mixture containing 81% of the (1S,1'R) diastereoisomer and 19% of 3 other diastereomers was obtained.. The addition of 60 parts by weight of I

to  
a standard perfuming composition imparted to the latter a powerful musky aspect,  
which is also smooth and fresh, together with a green, pear's peel connotation and a Galbanum undertone.

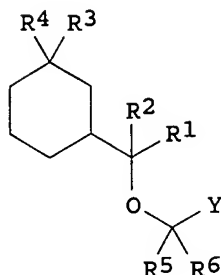
L8 ANSWER 3 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Methods for the production of novel alicyclic esters having a musky smell  
AN 2003:796643 CAPLUS  
DN 139:307907  
TI Methods for the production of novel alicyclic esters having a musky smell  
IN Eh, Marcus  
PA Symrise GmbH & Co. KG, Germany  
SO PCT Int. Appl., 45 pp.  
CODEN: PIXXD2

DT Patent  
LA German

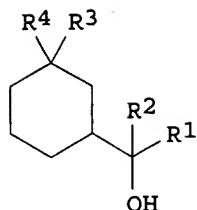
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2003082799	A2	20031009	WO 2003-EP3294	20030329
	WO 2003082799	A3	20031231		
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	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
				DE 2002-10214675	A 20020403
	DE 10214675	A1	20031016	DE 2002-10214675	20020403
	BR 2003004218	A	20040727	BR 2003-4218	20030329
				DE 2002-10214675	A 20020403
				WO 2003-EP3294	W 20030329
	EP 1492759	A2	20050105	EP 2003-732270	20030329
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK				

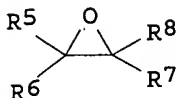
OS CASREACT 139:307907; MARPAT 139:307907  
GI



I



II



III

AB The invention relates to novel alicyclic esters I [R1 = Me; R2, R4 = H; R3 = H, Me; R5, R6 = H, Me; Y = CR7R8OC(:O)R9; R7, R8 = H, Me; R9 = C1-5-alkyl, C2-5-alkylene; or R1, R2 = Me, Et; R3, R4 = H, Me; R5R6 = O; Y = CR7R8OC(:O)R9; or R1, R2 = Me, Et; R4, R5, R6, R7 = H, Me; Y = CR7R8OC(:O)R9], methods for their production, for their use as odorous substances for perfumed products and for odorous substance mixts. containing the inventive compds. The procedure for the preparation of I is characterized by reaction of cyclohexylalkanols II with carboxylic acids [R9CO2CR7R8CO2H, R9CO2H or XCR7R8CO2H (X = OH, halogen)] anhydrides [(R9CO2)2O or (XCR7R8CO2)2O], or epoxides, III. Thus, I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2Cet] was prepared from 1-cyclohexylethanol via reaction with isobutylene oxide in cyclohexane containing BF3·OEt2, followed by reaction with (EtCO2)2O containing Et3N in the presence of catalytic DMAP. The odor of I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2Cet] was characterized (perceptible rose bloom note).

L8 ANSWER 4 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN

TI Preparation of cycloalkanecarboxylic acid derivatives for use as fragrances with musk characteristics

AN 2002:925306 CAPLUS

DN 138:4714

TI Preparation of cycloalkanecarboxylic acid derivatives for use as fragrances with musk characteristics

IN Kraft, Philip; Cadalbert, Riccardo

PA Givaudan SA, Switz.

SO Eur. Pat. Appl., 13 pp.

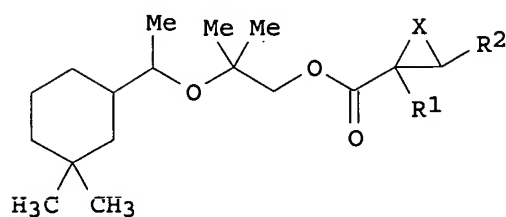
CODEN: EPXXDW

DT Patent

LA English

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	EP 1262474	A1	20021204	EP 2001-113377	20010601
	R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR				
	WO 2002096852	A1	20021205	WO 2002-CH282	20020530
	W: AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SD, SE, SG, SI, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VN, YU, ZA, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM				
	RW: GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG				
	EP 1392640	A1	20040303	EP 2001-113377	A 20010601
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				EP 2002-727147	20020530
				EP 2001-113377	A 20010601
				WO 2002-CH282	W 20020530
	BR 2002010098	A	20040413	BR 2002-10098	20020530
				EP 2001-113377	A 20010601
				WO 2002-CH282	W 20020530
	JP 2004535412	T2	20041125	JP 2003-500032	20020530
				EP 2001-113377	A 20010601
				WO 2002-CH282	W 20020530
	US 2004234568	A1	20041125	US 2004-478626	20040614
				EP 2001-113377	A 20010601
				WO 2002-CH282	W 20020530
OS	MARPAT 138:4714				
GI					



I

AB Odorant cycloalkanecarboxylic acid esters, such as I [R<sub>1</sub>, R<sub>2</sub> = H, Me; X = (CH<sub>2</sub>)<sub>n</sub>; n = 1, 2, 3; n plus number of carbon atoms in R<sub>1</sub> and R<sub>2</sub> is less than 5], with musk odor were prepared for use as ingredients in fine-fragrances and functional perfumery comps. Thus, I (R<sub>1</sub> = R<sub>2</sub> = H, X = CH<sub>2</sub>) was prepared by O-alkylation of α,3,3-trimethylcyclohexylmethanol with isobutylene oxide using EtAlCl<sub>2</sub> in toluene, and subsequent esterification of the resulting alc. with cyclopropanecarboxylic acid using DMAP and DCC in CH<sub>2</sub>Cl<sub>2</sub>. Fragrance formulations of the prepd cycloalkanecarboxylic acid esters were presented.

RE.CNT 3 THERE ARE 3 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 5 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN

TI A method for the calculation of odor character from molecular structure

AN 2002:610108 CAPLUS

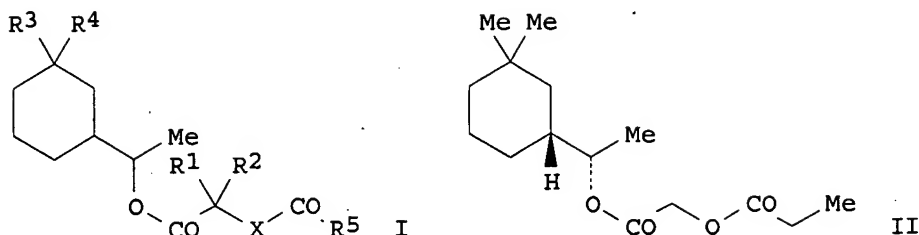
DN 139:3137  
 TI A method for the calculation of **odor** character from molecular structure  
 AU Turin, L.  
 CS Department of Physiology, University College London, London, WC1E 6BT, UK  
 SO Journal of Theoretical Biology (2002), 216(3), 367-385  
 CODEN: JTBIAP; ISSN: 0022-5193  
 PB Elsevier Science Ltd.  
 DT Journal  
 LA English  
 AB The relationship between mol. structure and **odor** character is one of the most complex structure-activity problems in biol. Despite over a century of effort, it remains unsolved, and synthesis of new odorants still proceeds largely by trial and error. In previous work, I have argued that the reason for this failure lies in a mistaken assumption, namely that mol. shape dets. **odor** character. Instead, I have taken up and extended an old idea (Dyson, 1938) according to which vertebrate olfactory receptors detect odorants by their mol. vibrations. I propose that the detection mechanism is inelastic electron tunnelling. If this is correct, there should be a correlation between the tunnelling vibrational spectra of odorants and their **odor** character. Here, using semi-empirical quantum chemical methods and a simple calcn. method for tunnelling mode intensities, I calculate the spectra of structurally diverse odorants belonging to various **odor** categories. With few exceptions, the calculated spectra of bitter almonds, musks, ambers, woods, sandalwoods and violets strongly correlate with **odor** character.  
 RE.CNT 28 THERE ARE 28 CITED REFERENCES AVAILABLE FOR THIS RECORD  
 ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 6 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
 TI Preparation of esters with musky **odor** and their use in perfumery  
 AN 2000:175782 CAPLUS  
 DN 132:194523  
 TI Preparation of esters with musky **odor** and their use in perfumery  
 IN Williams, Alvin S.  
 PA Firmenich S. A., Switz.  
 SO PCT Int. Appl., 24 pp.  
 CODEN: PIXXD2  
 DT Patent  
 LA English  
 FAN.CNT 1

PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
WO 2000014051	A1	20000316	WO 1999-IB1469	19990825
W: CA, CN, IL, IN, JP, US				
RW: AT, BE, CH, CY, DE, DK, ES, FI, FR, GB, GR, IE, IT, LU, MC, NL, PT, SE				
CA 2309449	AA	20000316	CH 1998-1841	A 19980909
			CA 1999-2309449	19990825
			CH 1998-1841	A 19980909
			WO 1999-IB1469	W 19990825
EP 1047660	A1	20001102	EP 1999-936930	19990825
EP 1047660	B1	20030129		
R: AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, FI				
			CH 1998-1841	A 19980909
			WO 1999-IB1469	W 19990825
JP 2002524438	T2	20020806	JP 2000-568811	19990825
			CH 1998-1841	A 19980909
			WO 1999-IB1469	W 19990825
AT 231834	E	20030215	AT 1999-936930	19990825
			CH 1998-1841	A 19980909
			WO 1999-IB1469	W 19990825
ES 2192067	T3	20030916	ES 1999-936930	19990825
			CH 1998-1841	A 19980909

US 6384269	B1	20020507	US 2000-553376	20000420
			WO 1999-IB1469	A1 19990825
US 38659	E	20041123	US 2002-326341	20021223
			CH 1998-1841	A 19980909
			WO 1999-IB1469	A1 19990825
			US 2000-553376	A5 20000420

OS MARPAT 132:194523  
GI



AB Cyclohexanemethanol esters I [R1, R2, R3, R4, R6, R7 = H, Me, Et; R5 = alkyl, alkoxy, alkenyl, acyloxy; X = CR6R7], which have a musky odor, were prepared for use as perfume additives. Thus, ester II was prepared with 79.5% yield by esterification of (+)-[S-(R\*,S\*)]- $\alpha$ ,3,3-trimethylcyclohexanemethanol with EtCO2CH2COCl. Perfuming compns. of the prepared compds. were presented.

RE.CNT 1 THERE ARE 1 CITED REFERENCES AVAILABLE FOR THIS RECORD  
ALL CITATIONS AVAILABLE IN THE RE FORMAT

L8 ANSWER 7 OF 7 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Synthesis of ethylene glycol mixed esters of alkanolic and cyclohexanoic acids  
AN 1968:77820 CAPLUS  
DN 68:77820  
TI Synthesis of ethylene glycol mixed esters of alkanolic and cyclohexanoic acids  
AU Zeinalov, B. K.; Radzhalov, D. T.; Dzhaifarova, A. A.; Muradyan, N. V.  
SO Azerbaidzhanskii Khimicheskii Zhurnal (1967), (2), 35-7  
CODEN: AZKZAU; ISSN: 0005-2531  
DT Journal  
LA Russian  
GI For diagram(s), see printed CA Issue.  
AB Acyloxyethyl esters (I and II) (tabulated) of cyclohexane- and 1,1-methylcyclohexanecarboxylic acids (sic) were prepared in 60-70% yield by heating a stirred mixture of 0.5 mole of the resp. alkanolic acid, 0.5 mole  $\beta$ -hydroxyethyl cyclohexanecarboxylate (II) or  $\beta$ -hydroxyethyl 1-methyl-1-cyclohexanecarboxylate (IV) 1% by weight of H2SO4, and 150 ml. dry C6H6 at 85-90° for 3-8 hrs., with a H2O aspirator in operation, distilling C6H6 from the neutralized, washed solution, and vacuum distilling the residue. The products were oily liqs. with pleasant ester odors and had good solubilizing and plasticizing properties. Poly(vinyl chloride) plasticized with these esters was distinguished by high elasticity. III, b1.5 110-12°, d20 1.0668, n20D 1.4680, and IV, b1 105-6°, d20 1.0436, n20D 1.4665, were prepared by condensing the resp. acids with ClCH2CH2OH in the presence of 40% aqueous alkali. [TABLE OMITTED]

=> file reg  
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FULL ESTIMATED COST

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ENTRY	SESSION
30.78	196.82



DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE ENTRY	TOTAL SESSION
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Property values tagged with IC are from the ZIC/VINITI data file provided by InfoChem.

STRUCTURE FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5  
 DICTIONARY FILE UPDATES: 17 JUL 2005 HIGHEST RN 855596-49-5

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TSCA INFORMATION NOW CURRENT THROUGH JANUARY 18, 2005

Please note that search-term pricing does apply when conducting SmartSELECT searches.

```
*****
*
* The CA roles and document type information have been removed from *
* the IDE default display format and the ED field has been added, *
* effective March 20, 2005. A new display format, IDERL, is now *
* available and contains the CA role and document type information. *
*
*****
```

Structure search iteration limits have been increased. See HELP SLIMITS for details.

Experimental and calculated property data are now available. For more information enter HELP PROP at an arrow prompt in the file or refer to the file summary sheet on the web at:  
<http://www.cas.org/ONLINE/DBSS/registryss.html>

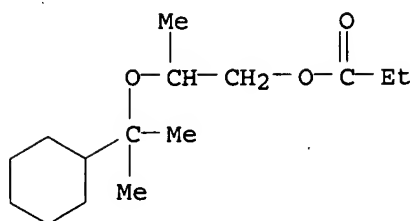
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=> e 1-Propanol, 2-(1-cyclohexyl-1-methylethoxy)-, propanoate/cn
E1      1      1-PROPANOL, 2-(1-BUTOXYETHOXY)-, (2R)-/CN
E2      1      1-PROPANOL, 2-(1-CYCLOHEXYL-1-METHYLETHOXY)-, ACETATE/CN
E3      1 --> 1-PROPANOL, 2-(1-CYCLOHEXYL-1-METHYLETHOXY)-, PROPANOATE/CN
E4      1      1-PROPANOL, 2-(1-CYCLOHEXYLETHOXY)-2-METHYL-/CN
E5      1      1-PROPANOL, 2-(1-CYCLOHEXYLETHOXY)-2-METHYL-, ACETATE/CN
E6      1      1-PROPANOL, 2-(1-CYCLOHEXYLETHOXY)-2-METHYL-, PROPANOATE/CN
E7      1      1-PROPANOL, 2-(1-CYCLOHEXYLPROPOXY)-2-METHYL-, ACETATE/CN
E8      1      1-PROPANOL, 2-(1-CYCLOHEXYLPROPOXY)-2-METHYL-, PROPANOATE/CN
E9      1      1-PROPANOL, 2-(1-ETHOXYETHOXY)-/CN
E10     1      1-PROPANOL, 2-(1-ETHOXYETHOXY)-, (2R)-/CN
E11     1      1-PROPANOL, 2-(1-ETHOXYETHOXY)-, (2R)-(PARTIAL)-/CN
E12     1      1-PROPANOL, 2-(1-ETHOXYETHOXY)-, (2S)-/CN
```

```
=> e3
L9      1 "1-PROPANOL, 2-(1-CYCLOHEXYL-1-METHYLETHOXY)-, PROPANOATE"/CN
```

```
=> d 19
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```
L9      ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN
RN      610769-92-1 REGISTRY
ED      Entered STN: 30 Oct 2003
CN      1-Propanol, 2-(1-cyclohexyl-1-methylethoxy)-, propanoate (9CI)
        (CA INDEX NAME)
FS      3D CONCORD
MF      C15 H28 O3
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SR CA  
LC STN Files: CA, CAPLUS, CASREACT



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

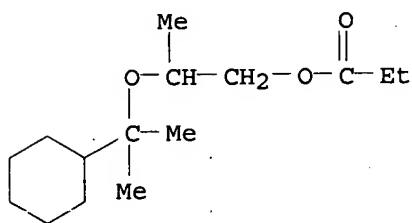
1 REFERENCES IN FILE CA (1907 TO DATE)  
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> search 19 exact full

L10 1 "1-PROPANOL, 2-(1-CYCLOHEXYL-1-METHYLETHOXY)-, PROPANOATE"/CN

=> d 110

L10 ANSWER 1 OF 1 REGISTRY COPYRIGHT 2005 ACS on STN  
RN 610769-92-1 REGISTRY  
ED Entered STN: 30 Oct 2003  
CN 1-Propanol, 2-(1-cyclohexyl-1-methylethoxy)-, propanoate (9CI)  
(CA INDEX NAME)  
FS 3D CONCORD  
MF C15 H28 O3  
SR CA  
LC STN Files: CA, CAPLUS, CASREACT



\*\*PROPERTY DATA AVAILABLE IN THE 'PROP' FORMAT\*\*

1 REFERENCES IN FILE CA (1907 TO DATE)  
1 REFERENCES IN FILE CAPLUS (1907 TO DATE)

=> file caplus

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
13.31	210.13

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
0.00	-5.11

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FILE COVERS 1907 - 19 Jul 2005 VOL 143 ISS 4  
FILE LAST UPDATED: 18 Jul 2005 (20050718/ED)

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This file contains CAS Registry Numbers for easy and accurate substance identification.

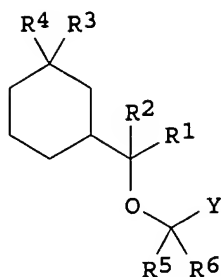
=> l10

L11 1 L10

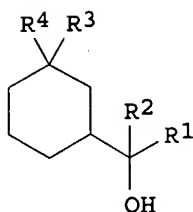
=> d l11 ti fbib abs

L11 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Methods for the production of novel alicyclic esters having a musky smell  
AN 2003:796643 CAPLUS  
DN 139:307907  
TI Methods for the production of novel alicyclic esters having a musky smell  
IN Eh, Marcus  
PA Symrise GmbH & Co. KG, Germany  
SO PCT Int. Appl., 45 pp.  
CODEN: PIXXD2  
DT Patent  
LA German  
FAN.CNT 1

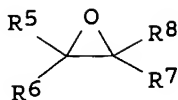
	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	WO 2003082799	A2	20031009	WO 2003-EP3294	20030329
	WO 2003082799	A3	20031231		
	W:	AE, AG, AL, AM, AT, AU, AZ, BA, BB, BG, BR, BY, BZ, CA, CH, CN, CO, CR, CU, CZ, DE, DK, DM, DZ, EC, EE, ES, FI, GB, GD, GE, GH, GM, HR, HU, ID, IL, IN, IS, JP, KE, KG, KP, KR, KZ, LC, LK, LR, LS, LT, LU, LV, MA, MD, MG, MK, MN, MW, MX, MZ, NO, NZ, OM, PH, PL, PT, RO, RU, SC, SD, SE, SG, SK, SL, TJ, TM, TN, TR, TT, TZ, UA, UG, US, UZ, VC, VN, YU, ZA, ZM, ZW			
	RW:	GH, GM, KE, LS, MW, MZ, SD, SL, SZ, TZ, UG, ZM, ZW, AM, AZ, BY, KG, KZ, MD, RU, TJ, TM, AT, BE, BG, CH, CY, CZ, DE, DK, EE, ES, FI, FR, GB, GR, HU, IE, IT, LU, MC, NL, PT, RO, SE, SI, SK, TR, BF, BJ, CF, CG, CI, CM, GA, GN, GQ, GW, ML, MR, NE, SN, TD, TG			
	DE 10214675	A1	20031016	DE 2002-10214675	20020403
	BR 2003004218	A	20040727	BR 2003-4218	20030329
				DE 2002-10214675	A 20020403
				WO 2003-EP3294	W 20030329
	EP 1492759	A2	20050105	EP 2003-732270	20030329
	R:	AT, BE, CH, DE, DK, ES, FR, GB, GR, IT, LI, LU, NL, SE, MC, PT, IE, SI, LT, LV, FI, RO, MK, CY, AL, TR, BG, CZ, EE, HU, SK			
				DE 2002-10214675	A 20020403
				WO 2003-EP3294	W 20030329
OS	CASREACT 139:307907; MARPAT 139:307907				
GI					



I



II



III

AB The invention relates to novel alicyclic esters I [R1 = Me; R2, R4 = H; R3 = H, Me; R5, R6 = H, Me; Y = CR7R8OC(:O)R9; R7, R8 = H, Me; R9 = C1-5-alkyl, C2-5-alkylene; or R1, R2 = Me, Et; R3, R4 = H, Me; R5R6 = O; Y = CR7R8OC(:O)R9; or R1, R2 = Me, Et; R4, R5, R6, R7 = H, Me; Y = CR7R8OC(:O)R9], methods for their production, for their use as odorous substances for perfumed products and for odorous substance mixts. containing the inventive compds. The procedure for the preparation of I is characterized by reaction of cyclohexylalkanols II with carboxylic acids [R9CO2CR7R8CO2H, R9CO2H or XCR7R8CO2H (X = OH, halogen)] anhydrides [(R9CO2)2O or (XCR7R8CO2)2O], or epoxides, III. Thus, I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2Cet] was prepared from 1-cyclohexylethanol via reaction with isobutylene oxide in cyclohexane containing BF3·OEt2, followed by reaction with (EtCO2)2O containing Et3N in the presence of catalytic DMAP. The odor of I [R1 = Me, R2 - R4 = H, R6 = CHMe2, Y = O2Cet] was characterized (perceptible rose bloom note).

=> d his

(FILE 'HOME' ENTERED AT 06:25:36 ON 19 JUL 2005)

FILE 'REGISTRY' ENTERED AT 06:27:51 ON 19 JUL 2005

L1 STRUCTURE UPLOADED  
L2 50 SEARCH L1 SSS SAM  
L3 STRUCTURE UPLOADED  
L4 0 SEARCH L3 SSS SAM  
L5 74 SEARCH L3 SSS FULL  
SAVE TEMP L5 RAWFINDS/A

FILE 'CAPLUS' ENTERED AT 06:33:54 ON 19 JUL 2005

L6 21 L5  
SAVE TEMP ODORREFS/A L6  
L7 69532 ODOR  
L8 7 L6 AND L7

FILE 'REGISTRY' ENTERED AT 06:44:54 ON 19 JUL 2005

E 1-PROPANOL, 2-(1-CYCLOHEXYL-1-METHYLETHOXY)-, PROPANOATE/CN

L9 1 E3

L10 1 SEARCH L9 EXACT FULL

FILE 'CAPLUS' ENTERED AT 06:45:43 ON 19 JUL 2005

L11 1 L10

=> 16 not 18

L12 14 L6 NOT L8

=> d 112 10-14 ti fbib abs

L12 ANSWER 10 OF 14 CAPLUS COPYRIGHT 2005 ACS on STN

TI Yellow dyes, ink compositions containing the same, and thermal transfer recording sheets using the same, giving high-density stable images resistant to heat, light, or humidity

AN 1998:693003 CAPLUS

DN 130:14995

TI Yellow dyes, ink compositions containing the same, and thermal transfer recording sheets using the same, giving high-density stable images resistant to heat, light, or humidity

IN Ogiso, Akira; Shimokawa, Yasushi; Ito, Naoto

PA Mitsui Chemicals Inc., Japan

SO Jpn. Kokai Tokkyo Koho, 17 pp.

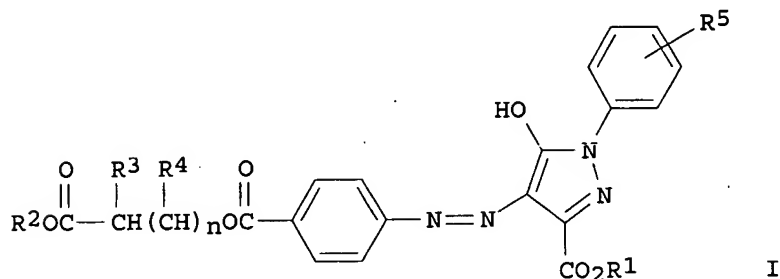
CODEN: JKXXAF

DT Patent

LA Japanese

FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	JP 10287056	A2	19981027	JP 1997-98915	19970416
	JP 3582954	B2	20041027		
				JP 1997-98915	19970416
OS	MARPAT 130:14995				
GI					



AB The title dyes have the general formula I, wherein R1, R2 = C1-10 (un)substituted alkyl, alkoxyalkyl; R3, R4 = H, Me; R5 = H, halogen, (un)substituted alkyl, alkoxy; n = 0, 1. An ink comprised I (R1 = R2 = Et; R3 = R5 = H; n = 0) 3, polybutyral resin 4.5, MEK 46.25, and toluene 46.25 parts.

L12 ANSWER 11 OF 14 CAPLUS COPYRIGHT 2005 ACS on STN

TI Branched bismethacrylates based on bis-GMA - a systematic route to low shrinkage composites

AN 1997:579234 CAPLUS

DN 127:221335

TI Branched bismethacrylates based on bis-GMA - a systematic route to low shrinkage composites

AU Holter, Dirk; Frey, Holger; Mulhaupt, Rolf  
 CS Institut für Makromolekulare Chemie und Freiburger  
 Materialforschungszentrum (FMF) der Albert-Ludwigs-Universität, Freiburg  
 i. Br., D-79104, Germany  
 SO Polymer Preprints (American Chemical Society, Division of Polymer  
 Chemistry) (1997), 38(2), 84-85  
 CODEN: ACPPAY; ISSN: 0032-3934  
 PB American Chemical Society, Division of Polymer Chemistry  
 DT Journal  
 LA English  
 AB Bismethacrylate monomers were prepared by esterification of bisphenol A  
 diglycidyl methacrylate (Bis-GMA) with different aliphatic and aromatic acids.  
 The relationship between mol. structure, viscosity, volume shrinkage and  
 mech. properties was studied. Variation of the side groups attached to  
 the Bis-GMA scaffold allows to decrease viscosity and volume shrinkage  
 concurrently. To obtain reasonable mech. properties Ph rings were  
 introduced into the side groups. 4-Phenylbutyric acid derivatized Bis-GMA  
 shows a viscosity of 51 Pas. The resp. composite with 50 vol% of a glass  
 filler reveals a volume shrinkage of 1.5% and an E-modulus of 6,300 MPa.

L12 ANSWER 12 OF 14 CAPLUS COPYRIGHT 2005 ACS on STN  
 TI Lipase-catalyzed resolution of 3-(aryloxy)-1,2-propanediol derivatives.  
 Towards an improved active site model of Pseudomonas cepacia lipase (Amano  
 PS)  
 AN 1995:663901 CAPLUS  
 DN 123:339233  
 TI Lipase-catalyzed resolution of 3-(aryloxy)-1,2-propanediol derivatives.  
 Towards an improved active site model of Pseudomonas cepacia lipase (Amano  
 PS)  
 AU Theil, Fritz; Lemke, Karin; Ballschuh, Sibylle; Kunath, Annamarie; Schick,  
 Hans  
 CS Inst. Angewandte Chemie Berlin-Adlershof e.V., Berlin, D-12484, Germany  
 SO Tetrahedron: Asymmetry (1995), 6(6), 1323-44  
 CODEN: TASYE3; ISSN: 0957-4166  
 PB Elsevier  
 DT Journal  
 LA English  
 OS CASREACT 123:339233  
 AB A variety of 3-(aryloxy)-1,2-propanediol derivs. with different  
 substituents on the aromatic ring or at the primary hydroxy group were used  
 as substrates for kinetic resolution by transesterification with vinyl  
 acetate catalyzed by lipase from Pseudomonas cepacia (Amano PS). Derivs.  
 with substituents in the para position of the aromatic ring were accepted as  
 substrates and resolved with high enantioselectivity. The corresponding  
 derivs. with substituents in the ortho position were much worse substrates  
 for lipase PS or even non-substrates if the substituent was sufficiently  
 space-filling, e.g., tert-Bu, Ph, benzyl or benzoyl. If the primary  
 hydroxy group was substituted by unbranched long-chain acyl residues, very  
 good substrates resulted. In contrast, derivs. with sterically crowded  
 residues at the primary hydroxy group (e.g., pivaloyl,  
 tert-butyldimethylsilyl, methanesulfonyl, p-toluenesulfonyl, or trityl)  
 were non-substrates for lipase PS.

L12 ANSWER 13 OF 14 CAPLUS COPYRIGHT 2005 ACS on STN  
 TI Preparation of 4-cycloalkyl-3-oxapentyl alkanoates as perfume fragrances  
 AN 1992:407513 CAPLUS  
 DN 117:7513  
 TI Preparation of 4-cycloalkyl-3-oxapentyl alkanoates as perfume fragrances  
 IN Giersch, Wolfgang Klaus; Schulte-Elte, Karl Heinrich  
 PA Firmenich S. A., Switz.  
 SO Eur. Pat. Appl., 15 pp.  
 CODEN: EPXXDW  
 DT Patent  
 LA French  
 FAN.CNT 1



FULL ESTIMATED COST	18.15	228.28
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-4.38	-9.49

SESSION WILL BE HELD FOR 60 MINUTES  
STN INTERNATIONAL SESSION SUSPENDED AT 06:48:34 ON 19 JUL 2005

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

\* \* \* \* \* RECONNECTED TO STN INTERNATIONAL \* \* \* \* \*  
SESSION RESUMED IN FILE 'CAPLUS' AT 07:14:36 ON 19 JUL 2005  
FILE 'CAPLUS' ENTERED AT 07:14:36 ON 19 JUL 2005  
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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	18.15	228.28
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-4.38	-9.49

=> d his

(FILE 'HOME' ENTERED AT 06:25:36 ON 19 JUL 2005)

FILE 'REGISTRY' ENTERED AT 06:27:51 ON 19 JUL 2005

L1 STRUCTURE UPLOADED  
L2 50 SEARCH L1 SSS SAM  
L3 STRUCTURE UPLOADED  
L4 0 SEARCH L3 SSS SAM  
L5 74 SEARCH L3 SSS FULL  
SAVE TEMP L5 RAWFINDS/A

FILE 'CAPLUS' ENTERED AT 06:33:54 ON 19 JUL 2005

L6 21 L5  
SAVE TEMP ODORREFS/A L6  
L7 69532 ODOR  
L8 7 L6 AND L7

FILE 'REGISTRY' ENTERED AT 06:44:54 ON 19 JUL 2005

E 1-PROPANOL, 2-(1-CYCLOHEXYL-1-METHYLETHOXY) -, PROPANOATE/CN  
L9 1 E3  
L10 1 SEARCH L9 EXACT FULL

FILE 'CAPLUS' ENTERED AT 06:45:43 ON 19 JUL 2005

L11 1 L10  
L12 14 L6 NOT L8

=> fragr?

L13 13878 FRAGR?

=> l6 and l13

L14 3 L6 AND L13



=> l14 not l8  
L15 1 L14 NOT L8

=> d l15 ti fbib abs

L15 ANSWER 1 OF 1 CAPLUS COPYRIGHT 2005 ACS on STN  
TI Preparation of 4-cycloalkyl-3-oxapentyl alkanoates as perfume  
fragrances  
AN 1992:407513 CAPLUS  
DN 117:7513  
TI Preparation of 4-cycloalkyl-3-oxapentyl alkanoates as perfume  
fragrances  
IN Giersch, Wolfgang Klaus; Schulte-Elte, Karl Heinrich  
PA Firmenich S. A., Switz.  
SO Eur. Pat. Appl., 15 pp.  
CODEN: EPXXDW  
DT Patent  
LA French  
FAN.CNT 1

	PATENT NO.	KIND	DATE	APPLICATION NO.	DATE
PI	EP 472966	A1	19920304	EP 1991-113240	19910807
	EP 472966	B1	19940928		
	R: CH, DE, FR, GB, LI, NL				
	US 5166412	A	19921124	CH 1990-2799	A 19900828
				US 1991-741027	19910806
				CH 1990-2799	A 19900828
	JP 06072952	A2	19940315	JP 1991-214881	19910827
	JP 2974834	B2	19991110		
				CH 1990-2799	A 19900828

OS MARPAT 117:7513  
AB RCHMeOCR1R2CR3R4O2CR5 (R = 3,3-dimethylcyclopentyl, -cyclohexyl; when R1 = R2 = H, R3 and/or R4 = Me; when R3 = R4 = H, R1 and/or R2 = Me; R5 = alkyl) were prepared Thus, 1-(3,3-dimethyl-1-cyclohexyl)-1-ethanone was ketalized by HOCH2CHMeOH and the dioxolane product reduced with Dibal to give RCHMeOCHMeCH2OR4 (I; R = 3,3-dimethylcyclohexyl) (II; R4 = H) and RCHMeOCH2CHMeOR4 (III; R same as I) (IV; R4 = H) as a mixture which was treated with EtCOCl to give II and IV (R4 = EtCO in each) as mixts. of diastereomers. Perfume formulations comprising title compds. are given.

=> d cost

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
4.68	16.37
0.72	2.70
1.89	173.88
18.55	43.02
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25.84	235.97

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION
-5.11	-10.22

CA SUBSCRIBER PRICE

IN FILE 'CAPLUS' AT 07:18:43 ON 19 JUL 2005

=> logoff hold

COST IN U.S. DOLLARS

SINCE FILE	TOTAL
ENTRY	SESSION
25.84	235.97

FULL ESTIMATED COST

DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)

SINCE FILE	TOTAL
ENTRY	SESSION

CA SUBSCRIBER PRICE

-5.11

-10.22

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 07:19:00 ON 19 JUL 2005

Connecting via Winsock to STN

Welcome to STN International! Enter x:x

LOGINID:SSSPTA1623PAZ

PASSWORD:

\* \* \* \* \* RECONNECTED TO STN INTERNATIONAL \* \* \* \* \*

SESSION RESUMED IN FILE 'CAPLUS' AT 08:01:54 ON 19 JUL 2005

FILE 'CAPLUS' ENTERED AT 08:01:54 ON 19 JUL 2005

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COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	25.84	235.97
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-5.11	-10.22

=> logoff hold

COST IN U.S. DOLLARS	SINCE FILE	TOTAL
	ENTRY	SESSION
FULL ESTIMATED COST	25.84	235.97
DISCOUNT AMOUNTS (FOR QUALIFYING ACCOUNTS)	SINCE FILE	TOTAL
	ENTRY	SESSION
CA SUBSCRIBER PRICE	-5.11	-10.22

SESSION WILL BE HELD FOR 60 MINUTES

STN INTERNATIONAL SESSION SUSPENDED AT 08:02:06 ON 19 JUL 2005